Nonlinear Transport and Current Fluctuation in an AB Ring with a Quantum Dot

JUNKO TAKAHASHI* and SHUICHI TASAKI

Department of Applied Physics, School of Science and Engineering, Waseda University, Tokyo 169-8555

Nonequilibrium steady states are explicitly constructed for a noninteracting electron model of an Aharonov-Bohm (AB) ring with a quantum dot (QD) with the aid of asymptotic fields. The Fano line shapes and AB oscillations are shown to strongly depend on the bias voltage. Current fluctuations are studied as well.

KEYWORDS: Fano effect, quantum dot, AB ring, current fluctuation, nonequilibrium steady states

1. Introduction

For an electron transport in an Aharonov-Bohm (AB) ring with a quantum dot (QD), a coupling between the discrete states and continuum causes an asymmetry of the resonance peak in the conductance, the so-called Fano resonance. As well known, when one varies a magnetic field piercing into the AB ring, the conductance oscillates (AB oscillation). And the phase of the AB oscillation changes by π from one side of the Fano resonance to the other. It is then interesting to see how these features appear in a far-from-equilibrium regime.

Recently, the method of C*-algebra has been widely applied to rigorous investigations of nonequilibrium steady states (NESS) of infinitely extended quantum systems. 4-9 The method was originally developed to deal with equilibrium statistical mechanics of such systems. 10,11 Fröhlich, Merkli and Ueltschi⁹ have shown the existence of NESS and its characterization for junction systems, which consist of several free-fermionic reservoirs mutually interacting via bounded interactions. Their results cover the AB ring with a QD in case that the electrodes are free fermion systems, but explicit expressions of observables at NESS are not given.

In this article, exact NESS is constructed explicitly for a noninteracting-electron model of the AB ring with a QD. We then examine the bias voltage dependence of the Fano resonance shapes and the AB oscillation. Near the Fano peak, AB oscillation patterns are shown to be strongly affected by the bias voltage. This feature is shown to imply nonlinear conduction which is sensitive to the magnetic field. Although we consider a noninteracting system, we believe that the Coulomb repulsion does not affect the essential features of the bias-voltage-dependence of the AB oscillation. The current fluctuation is investigated as well and we find that the interference effect appears in the fluctuation even when it cannot be seen in the average current.

2. System and Nonequilibrium Steady State

The system consists of two-dimensional electrodes and a quantum dot interacting with each other, which is schematically shown in Fig.1.¹² Only the dynamics of electrons is considered and the electron-electron interaction is neglected. Then, its Hamiltonian is given by

 $H = H_L + H_R + H_D + H_T + H_{LR}$, where $H_{\lambda} = \int dk \omega_{k\lambda} a^{\dagger}_{k\sigma\lambda} a_{k\sigma\lambda}$ ($\lambda = L$ (R)) is the energy of the left (right) electrode and $a_{k\sigma\lambda}$ ($\lambda = L$ (R)) is the annihilation operator of an electron with wave number k and spin σ at the corresponding electrode. Throughout this article, the repeated spin index means the summation over it. The bias voltage V is taken into account as the difference of the one-particle energies: $\omega_{kR} = \omega_{kL} - eV$. The second term $H_D = \epsilon_g c^{\dagger}_{\sigma} c_{\sigma}$ stands for the energy of the QD, where ϵ_g is the dot level and c_{σ} is the annihilation operator of the electron with spin σ at the dot. The electron transfer through the QD is described by a tunneling Hamiltonian

 $H_T = \int dk \{ u_{kL} a_{k\sigma L}^{\dagger} c_{\sigma} + u_{kR} a_{k\sigma R}^{\dagger} c_{\sigma} + (\text{H.c.}) \}$ and the direct transfer by

 $H_{LR} = \iint dk dq \{W_{kq} e^{i\varphi} a^{\dagger}_{k\sigma L} a_{q\sigma R} + (\text{H.c.})\},$ where $W_{kq} = W u_{kL} u_{qR}$ is assumed. The phase φ is proportional to the magnetic flux ϕ piercing between the two paths: $\varphi = e\phi/(\hbar c)$ with c the velocity of light.

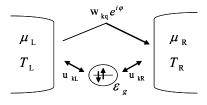


Fig. 1. Schematic view of an AB ring with a quantum dot

NESS can be simply characterized by in-coming fields $\beta_{k\sigma\lambda}$. They are solutions of $[H, \beta_{k\sigma\lambda}] = -\omega_{k\lambda}\beta_{k\sigma\lambda}$ with the boundary condition: $a_{k\sigma\lambda}(t) \exp(i\omega_{k\lambda}t/\hbar) \to \beta_{k\sigma\lambda}$ as $t \to -\infty$. When H does not have bound states, the original operators can be written as

$$a_{k\sigma\lambda} = \beta_{k\sigma\lambda} + \int dk' \left(\frac{u_{k\lambda} u_{k'\lambda}}{\omega_{k'\lambda} - \omega_{k\lambda} + i0} \frac{\chi_{k'\lambda} \beta_{k'\sigma\lambda}}{\Lambda(\omega_{k'\lambda})} + \frac{u_{k\lambda} u_{k'\bar{\lambda}}}{\omega_{k'\bar{\lambda}} - \omega_{k\lambda} + i0} \frac{\kappa_{k\bar{\lambda}} \beta_{k'\sigma\bar{\lambda}}}{\Lambda(\omega_{k'\bar{\lambda}})} \right)$$
(1)

$$c_{\sigma} = \int dk \left(\frac{u_{kL} \Omega_{kL} \beta_{k\sigma L}}{\Lambda(\omega_{kL})} + \frac{u_{kR} \Omega_{kR} \beta_{k\sigma R}}{\Lambda(\omega_{kR})} \right)$$
 (2)

where $\chi_{k\lambda} = 1 + M_{\lambda}(\omega_{k\lambda})[W^2(\omega_{k\lambda} - \epsilon_g) + 2W\cos\varphi]$, $\kappa_{k\lambda} = 1 + We^{i\varphi_{\lambda}}(\omega_{k\lambda} - \epsilon_g)$, $\Omega_{k\lambda} = 1 + We^{i\varphi_{\lambda}}M_{\lambda}(\omega_{k\lambda})$, $\Lambda(x) = \nu(x)(x - \epsilon_g) - \sum_{\lambda} M_{\lambda}(x) - 2W\cos\varphi M_L(x)M_R(x)$ $\nu(x) = 1 - W^2M_L(x)M_R(x)$, and

^{*}E-mail address: junchi@ruri.waseda.jp

$$M_{\lambda}(x) = \int d\tilde{k} \frac{u_{\tilde{k}\lambda}^2}{x - \omega_{\tilde{k}\lambda} + i0}.$$

In the above, $\bar{L} = R$, $\bar{R} = L$, $\varphi_L = -\varphi$ and $\varphi_R = \varphi$.

NESS is constructed as follows: (1) Initially, the tunneling interactions are turned off and the two electrodes are prepared to be in equilibrium with different temperatures $(T_L \text{ and } T_R)$ and chemical potentials $(\mu_L \text{ and } \mu_R)$. The initial dot state can be taken arbitrarily. Let ρ_0 be the density matrix describing this initial state. (2) The tunneling interactions are turned on. Then, the whole state $\rho_t \equiv e^{-iHt/\hbar}\rho_0 e^{iHt/\hbar}$ begins to evolve towards a steady state ρ_+ . One can actually show that, when $t \to +\infty$, ρ_t approaches to ρ_+ in a weak sense, namely, $\lim_{t\to +\infty} \operatorname{tr}(A\rho_t) = \operatorname{tr}(A\rho_+)$ holds for every local observable (for detail, see Appendix).

The so-obtained NESS ρ_+ satisfies Wick's theorem and is characterized by the two-point function:

$$\langle \beta_{k\sigma\lambda}^{\dagger} \beta_{k'\sigma'\lambda'} \rangle = F_{\lambda}(\epsilon) \delta(k - k') \delta_{\sigma\sigma'} \delta_{\lambda\lambda'} \tag{3}$$

where $\langle ... \rangle \equiv \operatorname{tr}(...\rho_{+})$ stands for the NESS average and $F_{\lambda}(\epsilon)$ denotes the Fermi distribution function with temperature T_{λ} and chemical potential μ_{λ} ($\lambda = L$ or R).

Here, a remark is in order. In spite of the fact that the system is conservative (i.e., the evolution is governed by the Hamiltonian), the state evolution exhibits 'dissipation' in the weak sense. This behavior is due to the infinite extension of the system. Indeed, information about the initial state is 'dissipated' towards infinity (namely the Poincaré recurrence cycle is infinite) and cannot be retrieved if we consider the average values of only the local observables. Thus, for every local observable A, the long-term limit $\lim_{t\to+\infty}\operatorname{tr}(A\rho_t)$ exists. Note that the Keldysh Green's function method shares the same property when applied to the description of NESS. In reality, the electrodes are finite, but can be regarded as infinite systems compared with the mesoscopic subsystems (AB-ring and QD). And the experimentally observed nonequilibrium states are considered to be welldescribed by NESS such as ρ_+ .

Now we go back to the steady-state current. The operator J corresponding to the current from the left electrode is $J = -\frac{d}{dt} \int dk \; a^{\dagger}_{k\sigma L} a_{k\sigma L}$ and its average is

$$\langle J \rangle = \frac{e}{h} \int dk u_{kL} J_k^{\text{dot}} + \frac{e}{h} \iint dk dq W u_{kL} u_{qR} J_{kq}^{\text{arm}}$$
 (4)

where $J_k^{\text{dot}} = i \langle a_{k\sigma L}^{\dagger} c_{\sigma} \rangle + (\text{c.c.})$ is the current via the QD and $J_{kq}^{\text{arm}} = i e^{i\varphi} \langle a_{k\sigma L}^{\dagger} a_{q\sigma R} \rangle + (\text{c.c.})$ is the current via the direct tunneling. With the aid of eqs.(1), (2) and (3), the current flowing from the left electrode is given by

$$\langle J \rangle = \frac{e}{h} \int_{\mu_c}^{\infty} d\epsilon \ T(\epsilon) \{ F_L(\epsilon) - F_R(\epsilon) \}$$
 (5)

where $\mu_c = \max(\omega_{0L}, \omega_{0R})$ and $T(\epsilon)$ is the transmission probability:

$$T(\epsilon) = \frac{4\pi^2 u^4 D^2}{|\Lambda(\epsilon)|^2} \{ (W(\epsilon - \epsilon_g) + \cos\varphi)^2 + \sin^2\varphi \}.$$

The constant D denotes the density of states of the two-dimensional electrodes. For the sake of the simplicity, the tunneling matrix elements u_{kL} , u_{kR} are taken to be symmetric and independent of the wave number:

 $u_{kL} = u_{kR} = u$. We remind that the bias voltage is equal to the chemical potential difference: $\mu_L - \mu_R = eV$. In the following sections, we focus on the nonequilibrium properties at zero temperature: $T_L = T_R = 0$.

3. Transport

The dot-level-dependence of the current is presented in Fig.2 for various bias voltages. In the linear response regime $(V \simeq 0)$, when the dot level passes through the Fermi energy of the electrodes, the resonant tunneling occurs and the current exhibits a typical Fano peak. When the bias voltage increases, the max-to-min distance becomes larger. This can be understood as follows. The current peak is generated by the interference between electrons passing through two conduction paths and this takes place when the dot energy ϵ_g lies between the two chemical potentials. Since the chemical potential difference is proportional to the bias voltage, the max-to-min distance of the current increases with the bias voltage. A similar deformation of the Fano peak as a function of the bias voltage was reported by Zhu $et\ al.^{13}$

The AB oscillations at a fixed dot level are depicted in Fig.3 for various bias voltages. The pattern of oscillation changes when one of the chemical potentials passes through the dot level. In particular, an oscillation with half wavelength appears when $|eV| = 0.10 \times \epsilon_F$ with ϵ_F the Fermi energy at zero bias.

As is known, ¹⁴ the two-terminal conductance G satisfies the Onsager relation $G(\varphi) = G(-\varphi)$ and, thus, is an even function of φ . This implies that the conductance takes its extremum at $\varphi = 0$ or π . Ueda et~al. ¹⁵ pointed out that the conductance can take its maximum at $\varphi \neq 0, \pi$. It is interesting to note that these two cases are simultaneously observed in Fig.3. When the bias voltage increases, the maximum of the current at $\varphi = 0$ splits into two maxima, which move to $\varphi = \pm \pi$ in a continuous way. But the minimum of the current abruptly jumps from $\varphi = \pi$ to $\varphi = 0$ at the bias voltage where the oscillation with half wavelength appears. Similar AB oscillation patterns can be seen in the linear response regime by varying the dot level. ¹⁶

The strong bias-voltage-dependence of the AB oscillation suggests a possibility of a device sensitive to the magnetic field. Indeed, the current-bias-voltage charac-

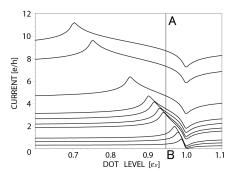


Fig. 2. Nonlinear current at zero temperature as a function of the gate voltage applied to the QD for different bias voltages eV from below eV/ϵ_F =0.01, 0.02, 0.03, 0.06, 0.07, 0.085, 0.10, 0.15, 0.25, 0.30. ϵ_F is the Fermi energy at eV = 0. The AB phase is fixed to φ = 0 and u^2DW = 0.1.

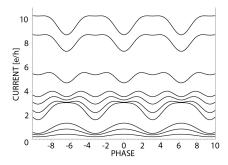


Fig. 3. Phase dependence of the current. The gate voltage is fixed to the A-B line shown in Fig.2. Other parameters are the same as those given in Fig.2.

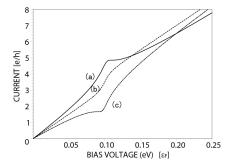


Fig. 4. Current versus bias voltage in units of e/h. The phase affects the current-bias-voltage characteristics: the phase (a) $\varphi = 0$ (the solid line), (b) $\varphi = \pi/2$ (the dashed line) and (c) $\varphi = \pi$ (the dotted line).

teristics at a fixed dot level depend on φ (Fig.4).

4. Current Fluctuations

Following Shimizu and Ueda, ¹⁷ we investigate the fluctuation of the charge $q(\tau) = \int_0^\tau dt J(t)$ leaving from the left electrode during the time interval τ , where J(t) is the current operator from the left electrode defined just before eq.(4). Its fluctuation

$$\Delta I \equiv \lim_{\tau \to \infty} \frac{\langle \{q(\tau) - \langle q(\tau) \rangle\}^2 \rangle}{\tau} \tag{6}$$

can be evaluated with the aid of Wick's theorem and the two-point function (3) and is given by

$$\Delta I = \left(\frac{e}{2\pi h}\right)^2 \int_{\mu_c}^{\infty} d\epsilon \left\{ T(\epsilon) \sum_{\lambda} F_{\lambda}(\epsilon) \left[1 - F_{\lambda}(\epsilon) \right] + T(\epsilon) \left[1 - T(\epsilon) \right] \left[F_L(\epsilon) - F_R(\epsilon) \right]^2 \right\}.$$
 (7)

The fluctuation ΔI consists of thermal and shot noises and, at zero temperature, only the shot noise survives.

Fig.5 shows a typical behavior of ΔI as a function of the dot level. One finds that ΔI has two dips, corresponding to the maximum and minimum of the current. Obviously, when the current is small, its fluctuation is small. And thus, the current minimum corresponds to the fluctuation dip. On the other hand, for one-dimensional perfect conductor, the current fluctuation is known to vanish¹⁷ because all the available states are used for the transport and there remains no room for fluctuations. In the same way, when the current takes its maximum, the dot level is used for the transport and is not available for

fluctuations. Then, the fluctuation is suppressed and the dip appears. Similar results are reported in connection with the photo-assisted shot noise. 18

When $\varphi = \pi/2$, the current is known to exhibit a symmetric Breit-Wigner-type resonance peak as for the system without direct tunneling between electrodes (the single dot system). This seems to indicate the absence of interference between electrons passing through two paths. However, it is not the case and the interference can be seen in the current fluctuation. Indeed, if the interference is absent, the fluctuation would be a sum of the fluctuation of the single dot system and constant fluctuation of the system without quantum dot. The curve (b) of the left figure of Fig.6 depicts the fluctuation of the present system at $\varphi = \pi/2$, clearly which is not a sum of a constant function and the fluctuation of the single dot system (the curve (b) of the right figure of Fig.6). The difference can be attributed to the interference between electrons passing through the two paths.

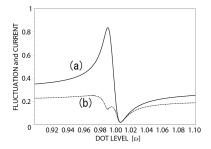


Fig. 5. The relationship between the current (a) and fluctuation (b) at low bias voltage range $(eV/\epsilon_F = 0.01)$, $\varphi = 0$, in units of (a) e/h and (b) $(e/2\pi h)^2$.

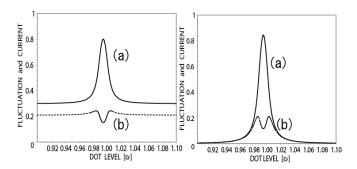


Fig. 6. The current (a) and fluctuation (b) with the present system at $\varphi = \pi/2$ (left) and with the single dot system (right) in units of (a) e/h and (b) $(e/2\pi h)^2$ at $eV/\epsilon_F = 0.01$ respectively.

5. Conclusions

In summary, exact NESS for the noninteracting model of an AB ring with a QD is explicitly constructed and bias-voltage-dependences of the Fano resonance and the AB oscillation are investigated. The AB oscillation patterns are shown to strongly depend on the bias voltage. This implies the nonlinear conduction sensitive to the magnetic field. The current fluctuations are studied as well and we show that, even when the interference cannot be seen in the average current, it contributes the current fluctuation.

Acknowledgment

The authors thank Professor T. Matsui for discussions. One of them (JT) is grateful to Professor S. Tarucha for fruitful discussions and valuable comments. This work is partially supported by a Grant-in-Aid for Scientific Research of Priority Areas "Control of Molecules in Intense Laser Fields" and the 21st Century COE Program at Waseda University "Holistic Research and Education Center for Physics of Self-organization Systems" both from the Ministry of Education, Culture, Sports, Science and Technology of Japan.

Appendix: C*-algebraic approach to NESS

In this approach, 10 one starts from a set of all (spatially) local observables, which form an algebra called the C*-algebra. Remind that, if the interaction range is finite, the commutator i[H,A] of the total Hamiltonian H and a local observable A is a local operator. Thus, both sides of the Heisenberg equation of motion $\hbar \frac{\partial A}{\partial t} = i[H,A]$ are local and the time evolution A(t) can be treated within the framework of local observables.

Physical states ρ are introduced by specifying average values $\langle A \rangle_{\rho}$ of all local observables A (i.e., the states are (positive) linear functionals over the algebra of local observables). Then, the long-term behavior of the system is naturally studied through the behaviors of the averages $\langle A \rangle_{\rho(t)}$. Then, for various systems, one can show the existence of the weak limits ρ_{\pm} defined by $\lim_{t\to\pm\infty}\langle A \rangle_{\rho(t)}=\langle A \rangle_{\rho_{\pm}}$ for any observable A. In this way, one can avoid the explicit use of infinite quantities such as the total Hamiltonian and total particle numbers, which are mathematically ill-defined, and one may realize the unidirectional state evolution in the weak sense.

As mentioned in the text, our model is a typical example of the junction systems studied by Frölich *et al.*⁹ Since they rigorously show the existence of NESS and give its mathematical characterizations, we give a heuristic proof of the existence of NESS and its characterization presented in the text.

Let $H_0 \equiv H_L + H_R$ and define the Møller operator $\Omega = \lim_{t \to -\infty} e^{iHt/\hbar} e^{-iH_0t/\hbar} = \lim_{t \to +\infty} e^{-iHt/\hbar} e^{iH_0t/\hbar}$, then the incoming field operators are expressed as $\beta_{k\sigma\lambda} = \lim_{t \to -\infty} e^{iHt/\hbar} e^{-iH_0t/\hbar} a_{k\sigma\lambda} e^{iH_0t/\hbar} e^{-iHt/\hbar} = \Omega a_{k\sigma\lambda} \Omega^{\dagger}$. Suppose that the total Hamiltonian H does not admit bound states, then, since Ω intertwines H with H_0 : $H\Omega = \Omega H_0$ and H_0 has bound states represented by c_σ , Ω should annihilate the states involving dot electrons and, thus, the product $\Pi \equiv \Omega^{\dagger}\Omega$ is a projection onto the states consisting only the electrons represented by $a_{k\sigma\lambda}$ (see the arguments in a standard text book on scattering theory¹⁹). Therefore, one has $\Pi = (1 - c_{\uparrow}^{\dagger}c_{\uparrow})(1 - c_{\downarrow}^{\dagger}c_{\downarrow})$ and it commutes with $a_{k\sigma\lambda}$. On the other hand, remind that the initial state can be written as $\rho_0 \propto \rho_c \exp\left[-\sum_{\lambda}(H_{\lambda} - \mu_{\lambda}N_{\lambda})/T_{\lambda}\right]$ where N_{λ} is the total number operator of each electrode and ρ_c is the initial dot state. Since ρ_0

commutes with H_0 , the steady state is given by $\rho_+ = \lim_{t \to +\infty} e^{-iHt/\hbar} e^{iH_0t/\hbar} \rho_0 e^{-iH_0t/\hbar} e^{iHt/\hbar}$. This seems to imply $\rho_+ = \Omega \rho_0 \Omega^{\dagger}$, but the normalization changes as in the wave function renormalization, and one has $\rho_+ = \Omega \rho_0 \Omega^{\dagger}/\mathrm{tr}(\Omega \rho_0 \Omega^{\dagger}) = \Omega \rho_0 \Omega^{\dagger}/\langle \Pi \rangle_0$, where $\langle \cdots \rangle_0$ stands for $\mathrm{tr}(\cdots \rho_0)$. Hence, for example,

$$\begin{split} &\langle \beta_{k_1\sigma_1\lambda_1}^\dagger \beta_{k_2\sigma_2\lambda_2}^\dagger \beta_{k_3\sigma_3\lambda_3} \beta_{k_4\sigma_4\lambda_4} \rangle \\ &= \frac{\langle \Pi a_{k_1\sigma_1\lambda_1}^\dagger \Pi a_{k_2\sigma_2\lambda_2}^\dagger \Pi a_{k_3\sigma_3\lambda_3} \Pi a_{k_4\sigma_4\lambda_4} \Pi \rangle_0}{\langle \Pi \rangle_0} \\ &= \langle a_{k_1\sigma_1\lambda_1}^\dagger a_{k_2\sigma_2\lambda_2}^\dagger a_{k_3\sigma_3\lambda_3} a_{k_4\sigma_4\lambda_4} \Pi \rangle_0 / \langle \Pi \rangle_0 \\ &= \langle a_{k_1\sigma_1\lambda_1}^\dagger a_{k_2\sigma_2\lambda_2}^\dagger a_{k_3\sigma_3\lambda_3} a_{k_4\sigma_4\lambda_4} \rangle_0 \ . \end{split}$$

The last equality follows from the fact that $a_{k\sigma\lambda}$ and c_{σ} (hence, Π) are statistically independent at the state ρ_0 . Therefore, since the initial state ρ_0 satisfies Wick's theorem for $a_{k\sigma\lambda}$, the same is true for ρ_+ and $\beta_{k\sigma\lambda}$ and the two-point functions $\langle \beta_{k\sigma\lambda}^{\dagger} \beta_{k'\sigma'\lambda'} \rangle$ are given by (3).

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